```
CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH,
             GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR,
             LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM,
             PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN,
             TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW, AM, AZ, BY,
             KG, KZ, MD, RU
        RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, BG,
             CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC,
             NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ,
             GW, ML, MR, NE, SN, TD, TG
                                           US 2003-627991
                                                             20030728
                            20040513
    US 2004091422
                       Α1
PRAI US 2002-398653P
                            20020727
                       Ρ
                            20021118
    US 2002-427266P
                       Ρ
                            20021230
                       Ρ
    US 2002-437270P
    MARPAT 140:141702
os
     533935-35-2P 533935-36-3P 533935-37-4P
IT
     533935-38-5P 653572-06-6P 653572-07-7P
     653572-08-8P 653572-09-9P 653572-10-2P
     653572-11-3P 653572-12-4P 653572-13-5P
     653572-14-6P 653572-15-7P 653572-16-8P
     653572-17-9P 653572-18-0P 653572-19-1P
     653572-20-4P 653572-21-5P 653572-22-6P
    RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU
     (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES
     (Uses)
        (design and synthesis of aminophosphinic acid derivs. as renal
        dipeptidase inhibitors and antitumor agents)
RN
     533935-35-2 CAPLUS
     2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-
     3-phenyl-, (2Z)- (9CI) (CA INDEX NAME)
```

Double bond geometry as shown.

RN 533935-36-3 CAPLUS
CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]3-phenyl-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 533935-37-4 CAPLUS CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-

Double bond geometry as shown.

RN 533935-38-5 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(4-iodophenyl)-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 653572-06-6 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(4-fluorophenyl)-, (2Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 653572-07-7 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(4-fluorophenyl)-, (2E)- (9CI) (CA INDEX NAME)

RN 653572-08-8 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(4-bromophenyl)-, (2Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 653572-09-9 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(3-iodophenyl)-, (2Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 653572-10-2 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(3-iodophenyl)-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 653572-11-3 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(2-iodophenyl)-, (2Z)- (9CI) (CA INDEX NAME)

RN 653572-12-4 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(2-iodophenyl)-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 653572-13-5 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-[4-(trifluoromethyl)phenyl]-, (2Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 653572-14-6 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-[4-(trifluoromethyl)phenyl]-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 653572-15-7 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(3,4-dichlorophenyl)-, (2Z)- (9CI) (CA INDEX NAME)

RN 653572-16-8 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(3,4-dichlorophenyl)-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 653572-17-9 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(3-chloro-4-fluorophenyl)-, (2Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 653572-18-0 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(3-chloro-4-fluorophenyl)-, (2E)- (9CI) (CA INDEX NAME)

RN 653572-19-1 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(3-bromo-4-fluorophenyl)-, (2Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 653572-20-4 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(3-bromo-4-fluorophenyl)-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 653572-21-5 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-[4-(diethylamino)phenyl]-, (2Z)- (9CI) (CA INDEX NAME)

RN 653572-22-6 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-[4-(diethylamino)phenyl]-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

L4 ANSWER 2 OF 2 CAPLUS COPYRIGHT 2004 ACS on STN

Aminophosphinic acid derivs. were synthesized as potential inhibitors of renal dipeptidase (RDP), an enzyme over-expressed in benign and malignant colon tumors, and several compds. showed potent enzyme-inhibitory activity. In an example reaction, (E)-[BocNH(C6H11CH2)CH]P(O)(OMe)[CH2C(CO2Me):CHPh] was prepared in two steps from [BocNH(C6H11CH2)CH]P(O)(OMe)H and hydrolyzed to (E)-[NH2(C6H11CH2)CH]P(O)(OH)[CH2C(CO2H):CHPh].

AN 2003:114414 CAPLUS

DN 139:6950

TI Design, synthesis and evaluation of new RDP inhibitors

AU Gurulingappa, Hallur; Buckhaults, Phillip; Kumar, Srinivas K.; Kinzler, Kenneth W.; Vogelstein, Bert; Khan, Saeed R.

CS The Sidney Kimmel Comprehensive Cancer Center at Johns Hopkins, Baltimore, MD, 21231, USA

SO Tetrahedron Letters (2003), 44(9), 1871-1873 CODEN: TELEAY; ISSN: 0040-4039

PB Elsevier Science Ltd.

DT Journal

LA English

OS CASREACT 139:6950

IT 533935-35-2P 533935-36-3P 533935-37-4P 533935-38-5P

RL: BSU (Biological study, unclassified); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)

(RDP inhibition activity; stereoselective preparation and RDP inhibition activity of aminophosphinic acid derivs.)

RN 533935-35-2 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-phenyl-, (2Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 533935-36-3 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-phenyl-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 533935-37-4 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(4-iodophenyl)-, (2Z)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RN 533935-38-5 CAPLUS

CN 2-Propenoic acid, 2-[[(1-amino-2-cyclohexylethyl)hydroxyphosphinyl]methyl]-3-(4-iodophenyl)-, (2E)- (9CI) (CA INDEX NAME)

Double bond geometry as shown.

RE.CNT 14 THERE ARE 14 CITED REFERENCES AVAILABLE FOR THIS RECORD ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> file uspatall COST IN U.S. DOLLARS	SINCE FILE	TOTAL
FULL ESTIMATED COST	ENTRY 9.95	SESSION 165.58
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	-1.39	-1.39

FILE 'USPATFULL' ENTERED AT 16:16:53 ON 28 MAY 2004

CA INDEXING COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

FILE 'USPAT2' ENTERED AT 16:16:53 ON 28 MAY 2004
CA INDEXING COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

=> d his

(FILE 'HOME' ENTERED AT 16:15:37 ON 28 MAY 2004)

FILE 'REGISTRY' ENTERED AT 16:15:58 ON 28 MAY 2004

L1 STRUCTURE UPLOADED

L2 0 S L1

L3 21 S L1 FUL

FILE 'CAPLUS' ENTERED AT 16:16:30 ON 28 MAY 2004

L4 2 S L3

FILE 'USPATFULL, USPAT2' ENTERED AT 16:16:53 ON 28 MAY 2004

=> s 13

L5 0 L3

=> file uspatall

COST IN U.S. DOLLARS SINCE FILE TOTAL

ENTRY SESSION

FULL ESTIMATED COST 3.54 169.12

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS) SINCE FILE TOTAL

ENTRY SESSION

CA SUBSCRIBER PRICE 0.00 -1.39

FILE 'USPATFULL' ENTERED AT 16:17:17 ON 28 MAY 2004
CA INDEXING COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

FILE 'USPAT2' ENTERED AT 16:17:17 ON 28 MAY 2004
CA INDEXING COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

=> s 13

L6 0 L3